

Seminar über Theoretische Chemie

jeweils donnerstags, 15:15 Uhr, im Seminarraum 408, 4. OG, Geb. 30.44

- 03.11.11** **Dr. Michael Harding**
Towards highly accurate *ab initio* thermochemistry of larger systems: Benzene
- 10.11.11** **Dr. Cong Wang**
Analytical evaluations of unlinked exponential-correlated one-center three- and four-electron integrals
- 17.11.11** **Dr. Michiel van Setten** (INT, Campus Nord)
The *GW* method for quantum chemistry applications:
Theory, implementation, benchmarks
- 24.11.11^{*)}** **Dr. Alexander Schug** (Steinbuch Centre for Computing)
14:30 Uhr In-silico protein folding using coarse-grained models
- 08.12.11** **Tilman Bodenstein**
Spin-Bahn-Effekte in mehrkernigen Übergangsmetallkomplexen
- 19.01.12** **Jun.-Prof. Dr. Thomas Kühne** (Universität Mainz)
Quantum chemistry in a glass of water
- 26.01.12** **Anikó Udvarhelyi** (MPI für medizinische Forschung, Heidelberg)
Photoinduced proton-coupled electron transfer in BLUF photoreceptors.
A CASSCF study.
- 02.02.12** **Dr. Andrey Yachmenev**
A simple method to improve the accuracy of the RI approximation for
two-electron Coulomb integrals
- 09.02.12** **Anna Hehn**
Calculation of molecular spectra using Fock-space coupled-cluster theory

***) Bitte die Uhrzeit beachten.**